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INFORMATION DISCLOSURE STATEMENT BY APPLICANT

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Sheet 1 of 2

Complete if Known	
Application Number	10/825,186
Filing Date	April 16, 2004
First Named Inventor	Zhang et al.
Art Unit	~~1645~~ 1631
Examiner Name	~~To Be Assigned~~ Moran
Attorney Docket Number	57953/1221 (ZHA01-01)

U.S. PATENT DOCUMENTS

Examiner Initials*	Cite No.[1]	U.S. Patent Document Number - Kind Code[2] (if known)	Publication Date MM-DD-YYYY	Name of Patentee or Applicant of Cited Document	Pages, Columns, Lines, Where Relevant Passages or Relevant Figures Appear
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FOREIGN PATENT DOCUMENTS

Examiner Initials*	Cite No.[1]	Foreign Patent Document: Country Code[3] Number[4] Kind Code[5] (if known)	Publication Date MM-DD-YYYY	Name of Patentee or Applicant of Cited Document	Pages, Columns, Lines, Where Relevant Passages or Relevant Figures Appear	T[6]

OTHER PRIOR ART – NON PATENT LITERATURE DOCUMENTS

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/M.M./	1	Chen et al., "Fractionation of Peptide with Disulfide Bond for Quantum Mechanical Calculation of Interaction Energy with Molecules," *Journal of Chemical Physics* 120(2):839-844 (2004)	
/M.M./	2	Chen et al., "Theoretical Method for Full *ab initio* Calculation of DNA/RNA-Ligand Interaction Energy," *Journal of Chemical Physics* 120(24):11386-11391 (2004)	
/M.M./	3	Gao et al., "An Efficient Linear Scaling Method for ab initio Calculation of Electron Density of Proteins," *Chemical Physics Letters* 394:293-297 (2004)	

Examiner Signature	/Marjorie Moran/	Date Considered	03/18/2007

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INFORMATION DISCLOSURE STATEMENT BY APPLICANT

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Sheet	2	of	2

Complete if Known	
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Filing Date	April 16, 2004
First Named Inventor	Zhang et al.
Group Art Unit	~~1645~~ 1631
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/M.M./	4	Xiang et al., "Fully Quantum Mechanical Energy Optimization for Protein-Ligand Structure," *Journal of Computational Chemistry* 25(12):1431-1437 (2004)	
	5	Zhang et al., "Full *ab initio* Computation of Protein-Water Interaction Energies," *Journal of Theoretical and Computational Chemistry* 3(1):43-49 (2004)	
	6	Zhang et al., "Molecular Caps for Full Quantum Mechanical Computation of Peptide-Water Interaction Energy," *Journal of Computational Chemistry* 24(15):1846-1852 (2003)	
	7	Zhang et al., "Molecular Fractionation with Conjugate Caps for Full Quantum Mechanical Calculation of Protein-Molecule Interaction Energy," *Journal of Chemical Physics* 119(7):3599-3605 (2003)	
	8	Zhang et al., "New Advance in Computational Chemistry: Full Quantum Mechanical ab Initio Computation of Streptavidin-Biotin Interaction Energy," *J. Phys. Chem.* 107:12039-12041 (2003)	
	9	Zhang et al., "Quantum Mechanical Map for Protein-Ligand Binding with Application to β-Trypsin/Benzamidine Complex," *Journal of Chemical Physics* 120(3):1145-1148 (2004)	

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